Amendments To The Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

What is claimed is:

1. (Currently Amended) A compound of formula formual (I):

$$Ar^{1} - CHCH_{2}NHCR^{1}R^{2}(CH_{2})_{m} - O - (CH_{2})_{p}CR^{1a}R^{2a} - Ar^{2}$$

$$OH \qquad (I)$$

or a salt, solvate, or physiologically functional derivative thereof, wherein:

Ar¹ is a group selected from

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wherein R⁴ represents hydrogen, halogen, -(CH₂)_qOR⁷, -NR⁷C(O)R⁸, -NR⁷SO₂R⁸, -SO₂NR⁷R⁸, -NR⁷R⁸, -OC(O)R⁹ or OC(O)NR⁷R⁸, and R³ represents hydrogen, halogen or C₁₋₄ alkyl;

or R⁴ represents –NHR¹⁰ and R³ and –NHR¹⁰ together form a 5- or 6-membered heterocyclic ring;

R⁵ represents hydrogen, halogen, –OR⁷ or –NR⁷R⁸;

R⁶ represents hydrogen, haloC₁₋₄alkyl, -OR⁷, −NR⁷R⁸, -OC(O)R⁹ or OC(O)NR⁷R⁸;

 R^7 and R^8 each independently represents hydrogen or C_{1-4} alkyl, or in the groups $-NR^7R^8$, $-SO_2NR^7R^8$ and $-OC(O)NR^7R^8$, R^7 and R^8 independently represent hydrogen or C_{1-4} alkyl or together with the nitrogen atom to which they are attached form a 5-, 6- or 7- membered nitrogen-containing ring,

R⁹ represents an aryl (eg phenyl or naphthyl) group which may be unsubstituted or substituted by one or more substituents selected from halogen, C₁₋₄ alkyl, hydroxyl, C₁₋₄ alkoxy or halo C₁₋₄ alkyl; and

q is zero or an integer from 1 to 4;

Ar² is a group:

$$R^{12}$$
 or R^{12} R^{11}

wherein

 R^{11} is selected from hydrogen, $C_{1\text{-6}}$ alkyl, hydroxy, $C_{1\text{-6}}$ alkoxy, cyano, nitro, halo, $C_{1\text{-6}}$ haloalkyl, XCO_2R^{16} , $-XC(O)NR^{15}R^{16}$, $-XNR^{14}C(O)R^{15}$, $-XNR^{14}C(O)R^{15}R^{16}$, $-XNR^{14}C(O)R^{15}R^{16}$, $-XNR^{14}SO_2R^{15}$, $-XSO_2NR^{17}R^{18}$, XSR^{14} , $XSOR^{14}$, XSO_2R^{14} , $-XNR^{15}R^{16}$, $-XNR^{14}C(O)OR^{15}$, or $XNR^{14}SO_2NR^{15}R^{16}$, or R^{11} is selected from -X-aryl, -X-hetaryl, or -X-(aryloxy), each optionally substituted by 1 or 2 groups independently selected from hydroxy, $C_{1\text{-6}}$ alkoxy, halo, $C_{1\text{-6}}$ alkyl, $C_{1\text{-6}}$ haloalkyl, cyano, nitro, $CONR^{15}R^{16}$, $-NR^{14}C(O)R^{15}$, SR^{14} , SOR^{14} , $-SO_2R^{14}$, $-SO_2NR^{17}R^{18}$, $-CO_2R^{16}$, $-NR^{15}R^{16}$, or hetaryl optionally substituted by 1 or 2 groups independently selected from hydroxy, $C_{1\text{-6}}$ alkoxy, halo, $C_{1\text{-6}}$ alkyl, or $C_{1\text{-6}}$ haloalkyl;

X is $-(CH_2)_r$ – or C_{2-6} alkenylene;

r is an integer from 0 to 6, preferably 0 to 4;

 R^{14} and R^{15} are independently selected from hydrogen, C_{1-6} alkyl, C_{3-7} cycloalkyl, aryl, hetaryl, hetaryl(C_{1-6} alkyl)- and aryl(C_{1-6} alkyl)- and R^{14} and R^{15} are each independently optionally substituted by 1 or 2 groups independently selected from halo, C_{1-6} alkyl,

 C_{3-7} cycloalkyl, C_{1-6} alkoxy, C_{1-6} haloalkyl, -NHC(O)(C_{1-6} alkyl), -SO₂(C_{1-6} alkyl), -SO₂(aryl), -CO₂H, and -CO₂(C_{1-4} alkyl), -NH₂, -NH(C_{1-6} alkyl), aryl(C_{1-6} alkyl)-, aryl(C_{2-6} alkenyl)-,

aryl(C_{2-6} alkynyl)-, hetaryl(C_{1-6} alkyl)-, -NHSO $_2$ aryl, -NH(hetaryl C_{1-6} alkyl), -NHSO $_2$ hetaryl,

-NHSO₂(C₁₋₆alkyl), -NHC(O)aryl, or -NHC(O)hetaryl:

or R¹⁴ and R¹⁵, together with the nitrogen atom to which they are bonded, form a 5-, 6- or 7- membered nitrogen – containing ring;

or where R¹¹ is –XNR¹⁴C(O)NR¹⁵R¹⁶, R¹⁴ and R¹⁵ may, together with the – NC(O)N- portion of the group R¹ to which they are bonded, form a saturated or

unsaturated ring, preferably a 5, 6, or 7-membered ring, for example an imidazolidine-ring, such as imidazolidine-2,4-dione;

or where R¹¹ is –XNR¹⁴C(O)OR¹⁵, R¹⁴ and R¹⁵ may, together with the – NC(O)O- portion of the group R¹¹ to which they are bonded, form a saturated or unsaturated ring, preferably a 5-, 6-, or 7- membered ring, for example an exazolidine ring, such as exazolidine 2.4-dione:

R¹⁶ is selected from hydrogen, C₁₋₆alkyl and C₃₋₇ cycloalkyl;

or where R¹¹ is –XC(O)NR¹⁵R¹⁶ or –XNR¹⁴C(O)NR¹⁵R¹⁶, R¹⁵ and R¹⁶ may, together with the nitrogen to which they are bonded, form a 5-, 6-, or 7-membered nitrogen containing ring;

 R^{17} and R^{18} are independently selected from hydrogen, C_{1-6} alkyl, C_{3-7} cycloalkyl, aryl, hetaryl, hetaryl(C_{1-6} alkyl)- and aryl(C_{1-6} alkyl)-, or R^{17} and R^{18} , together with the nitrogen to which they are bonded, form a 5-, 6-, or 7- membered nitrogen containing ring;

and R¹⁷ and R¹⁸ are each optionally substituted by one or two groups independently selected from halo, C₁₋₆alkyl, and C₃₋₇cycloalkyl, C₁₋₆haloalkyl;

 R^{12} is selected from hydrogen, pyridine, C_{1-6} alkyl, C_{1-6} alkoxy, halo, aryl, aryl(C_{1-6} alkyl)-, C_{1-6} haloalkoxy, and C_{1-6} haloalkyl;

 R^{13} is selected from hydrogen, hydroxy, C_{1-6} alkyl, C_{1-6} alkoxy, halo, aryl, aryl(C_{1-6} alkyl)-, C_{1-6} haloalkoxy, and C_{1-6} haloalkyl;

 R^1 and R^2 are independently selected from hydrogen and C_{1-4} alkyl with the proviso that the total number of carbon atoms in R^1 and R^2 is not more than 4;

one of R^{1a} and R^{2a} is selected from hydrogen and C₁₋₄alkyl, and the other of R^{1a} and R^{2a} represents hydrogen or C₁₋₄alkyl;

m is an integer of from 1 to 3; n is an integer of from 1 to 4; and p is zero or an integer of from 1 to 3;

and ____ represents a single or double bond.

2. (Currently Amended) A compound of formula (I) as defined in claim 1, or a salt, solvate or physiologically functional derivative thereof, wherein except that:

R^{1a} and R^{2a} each represent hydrogen;

and in the group Ar¹, either:

 R^4 represents halogen, -(CH₂)_qOR⁷, -NR⁷C(O)R⁸, -NR⁷SO₂R⁸, -SO₂NR⁷R⁸, -NR⁷R⁸.

-OC(O)R⁹ or OC(O)NR⁷R⁸, and R³ represents hydrogen or C₁₋₄ alkyl; or

R⁴ represents –NHR¹⁰ and R³ and –NHR¹⁰ together form a 5- or 6- membered heterocyclic ring;

- 3. (Currently Amended) A compound of formula (I) according to either claim 1 or claim 2 wherein the group Ar¹ is selected from groups (a) and (b) as defined in claim 1.
- 4. (Currently Amended) A compound of formula (I) according to <u>claim 1</u> any of claims 1 to 3 wherein, in the group Ar^2 , R^{11} is selected from hydrogen, C_{1-4} alkyl, hydroxy, halo, -NR¹⁴C(O)NR¹⁵R¹⁶, -NR¹⁴SO₂R¹⁵ and XSO₂NR¹⁷R¹⁸ wherein R¹⁴ to R¹⁸ are as defined in claim 1.
- 5. (Currently Amended) A compound of formula (I) according to <u>claim 1</u> any of claims 1 to 3 wherein, in the group Ar^2 , R^{11} is selected from cyano, $CONR^{15}R^{16}$, SR^{14} , SOR^{14} and SO_2R^{14} , wherein R^{14} , R^{15} and R^{16} are as defined in claim 1.

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- 6. (Currently Amended) A compound of formula (I) according to <u>claim 1</u> any of claims 1 to 5 wherein R¹² and R¹³ each represent hydrogen.
- 7. (Currently Amended) A compound of formula (I) according to <u>claim 1</u> any of claims 1 to 3 wherein R^{11} represents hydrogen and R^{12} and R^{13} each represent halogen or C_{1-6} alkyl.
- 8. (Currently Amended) A compound of formula (I) according to <u>claim 1</u> any of claims 1 to 7 wherein R¹ and R² are both hydrogen.
- 9. (Currently Amended) A compound of formula (I) according to <u>claim 1</u> any of claims 1 to 8 wherein each of m and n is independently 1 or 2, and p is zero or 1.
- 10. (Currently Amended) A compound of formula (I) according to claim 1 selected from:
- $4-((1R)-2-\{[2-((3R)-3-\{[(2,6-Dichlorobenzyl)oxy]methyl\}-2,3-dihydro-1,4-dih$
- benzodioxin-6-yl)ethyl]amino}-1-hydroxyethyl)-2-(hydroxymethyl)phenol;
- $4-\{(1R)-2-[(2-\{(3R)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3R)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3R)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3R)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3R)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3R)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-(Benzyloxy)methyl-2-[(Benzyloxy)methyl$
- yl}ethyl)amino]-1-hydroxyethyl}-2-(hydroxymethyl)phenol;
- $4-\{(1R)-2-[(2-\{(3S)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3S)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3S)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3S)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3S)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3S)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3S)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3S)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(2-\{(3S)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-(1R)-2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-(Benzyloxy)methyl-2-[(Benzy$
- yl}ethyl)amino]-1-hydroxyethyl}-2-(hydroxymethyl)phenol;
- 2-(Hydroxymethyl)-4- $\{(1R)$ -1-hydroxy-2- $[(2-\{(3R)$ -3- $[(\Box yridine$ -3-
- ylmethoxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-yl}ethyl)amino]ethyl}phenol;
- $4-((1R)-2-\{[2-((3R)-3-\{[(6-Chloropyridin-3-yl)methoxy]methyl\}-2,3-dihydro-1,4-k-((1R)-2-\{[(3R)-3-\{[(6-Chloropyridin-3-yl)methoxy]methyl\}-2,3-dihydro-1,4-k-((1R)-2-\{[(6-Chloropyridin-3-yl)methoxy]methyl\}-2,3-dihydro-1,4-k-((1R)-2-\{[(6-Chloropyridin-3-yl)methoxy]methyl\}-2,3-dihydro-1,4-k-((1R)-2-\{[(6-Chloropyridin-3-yl)methoxy]methyl]-2,3-dihydro-1,4-k-((1R)-2-\{[(6-Chloropyridin-3-yl)methoxy]methyl]-2,3-dihydro-1,4-k-((1R)-2-\{[(6-Chloropyridin-3-yl)methoxy]methyl]-2,3-dihydro-1,4-k-((1R)-2-\{[(6-Chloropyridin-3-yl)methoxy]methyl]-2,3-dihydro-1,4-k-((1R)-2-\{[(6-Chloropyridin-3-yl)methoxy]methyl]-2,3-dihydro-1,4-k-((1R)-2-\{[(6-Chloropyridin-3-yl)methoxy]methyl]-2,3-dihydro-1,4-k-((1R)-2-\{[(6-Chloropyridin-3-yl)methoxy]methyl]-2,3-dihydro-1,4-k-((1R)-2$
- benzodioxin-6-yl)ethyl]amino}-1-hydroxyethyl)-2-(hydroxymethyl)phenol;
- 4-((1R)-2-{[2-((3R)-3-{[(2,6-Dichloropyridin-3-yl)methoxy]methyl}-2,3-dihydro-
- 1,4-benzodioxin-6-yl)ethyl]amino}-1-hydroxyethyl)-2-(hydroxymethyl)phenol;
- 4-{(1R)-2-[(2-{2-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-
- yl}ethyl)amino]-1-hydroxyethyl}-2-(hydroxymethyl)phenol;
- 4-((1*R*)-2-{[2-((3*R*)-3-{[(5-Bromopyridin-3-yl)methoxy]methyl}-2,3-dihydro-1,4-benzodioxin-6-yl)ethyl]amino}-1-hydroxyethyl)-2-(hydroxymethyl)phenol;

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3-[({(2R)-7-[2-({(2R)-2-Hydroxy-2-[4-hydroxy-3-
(hydroxymethyl)phenyl]ethyl}amino)ethyl]-2,3-dihydro-1,4-benzodioxin-2-
yl}methoxy)methyl]benzonitrile;
3-[({(2R)-7-[2-({(2R)-2-Hydroxy-2-[4-hydroxy-3-
(hydroxymethyl)phenyl]ethyl}amino)ethyl]-2,3-dihydro-1,4-benzodioxin-2-
yl}methoxy)methyl]benzamide;
4-[(1R)-2-({2-[(3R)-3-({[3-(Cyclopentylthio)benzyl]oxy}methyl)-2,3-dihydro-1,4-
benzodioxin-6-yl]ethyl]amino)-1-hydroxyethyl]-2-(hydroxymethyl)phenol;
4-[(1R)-2-({2-[(3R)-3-({[3-(Cyclopentylsulfonyl)benzyl]oxy}methyl)-2,3-dihydro-
1,4-benzodioxin-6-yl]ethyl}amino)-1-hydroxyethyl]-2-(hydroxymethyl)phenol;
2-(Hydroxymethyl)-4-{(1R)-1-hydroxy-2-[(2-{(3R)-3-[(5-[4-
(methylsulfinyl)phenyl] []yridine-3-yl}methoxy)methyl]-2,3-dihydro-1,4-
benzodioxin-6-yl}ethyl)amino]ethyl}phenol;
N-\{3-[(\{(2R)-7-[2-(\{(2R)-2-Hydroxy-2-[4-hydroxy-3-
(hydroxymethyl)phenyl]ethyl}amino)ethyl]-2,3-dihydro-1,4-benzodioxin-2-
yl}methoxy)methyl]phenyl}urea;
4-((1R)-2-{[2-((3R)-3-{[(4-Chlorobenzyl)oxy]methyl}-2,3-dihydro-1,4-
benzodioxin-6-yl)ethyl]amino}-1-hydroxyethyl)-2-(hydroxymethyl)phenol;
4-((1R)-2-{[2-((3R)-3-{[(4-Fluorobenzyl)oxy]methyl}-2,3-dihydro-1,4-
benzodioxin-6-yl)ethyl]amino}-1-hydroxyethyl)-2-(hydroxymethyl)phenol;
4-((1R)-2-{[2-((3R)-3-{[(3,5-Dimethylbenzyl)oxy]methyl}-2,3-dihydro-1,4-
benzodioxin-6-yl)ethyl]amino}-1-hydroxyethyl)-2-(hydroxymethyl)phenol;
2-(Hydroxymethyl)-4-{(1R)-1-hydroxy-2-[(2-{(3R)-3-[(1-phenylethoxy)methyl]-
2,3-dihydro-1,4-benzodioxin-6-yl}ethyl)amino]ethyl}phenol;
2-(Hydroxymethyl)-4-[(1R)-1-hydroxy-2-({2-[(3R)-3-({[3-
(methylsulfonyl)benzyl]oxy}methyl)-2,3-dihydro-1,4-benzodioxin-6-
yl]ethyl}amino)ethyl]phenol;
4-((1R)-2-{[2-((3R)-3-{[3-(2,6-Dichlorophenyl)propoxy]methyl}-2,3-dihydro-1,4-
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benzodioxin-6-yl)ethyl]amino}-1-hydroxyethyl)-2-(hydroxymethyl)phenol;
3-[({(2R)-7-[2-({(2R)-2-Hydroxy-2-[4-hydroxy-3-(hydroxymethyl)phenyl]ethyl}amino)ethyl]-2,3-dihydro-1,4-benzodioxin-2-yl}methoxy)methyl]benzenesulfonamide;

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6-{2-[(2-{(3*R*)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-yl}ethyl)amino]-1-hydroxyethyl}-2-(hydroxymethyl)□yridine-3-ol;
N-(5-{(1*R*)-2-[(2-{(3*R*)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-yl}ethyl)amino]-1-hydroxyethyl}-2-hydroxyphenyl)methanesulfonamide;
4-{(1*R*)-2-[(2-{(3*R*)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-yl}ethyl)amino]-1-hydroxyethyl}-2-fluorophenol;
4-{(1*R*)-2-[(2-{(3*R*)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-yl}ethyl)amino]-1-hydroxyethyl}-3-methylphenol;
(1*R*)-1-(4-Amino-3,5-dichlorophenyl)-2-[(2-{(3*R*)-3-[(benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-yl}ethyl)amino]ethanol;
5-{(1*R*)-2-[(2-{(3*R*)-3-[(Benzyloxy)methyl]-2,3-dihydro-1,4-benzodioxin-6-yl}ethyl)amino]-1-hydroxyethyl}-2-hydroxyphenylformamide;

or a salt, solvate or physiologically functional derivative thereof.

11. (Currently Amended) A method for the prophylaxis or treatment of a clinical condition in a mammal, such as a human, for which a selective β₂-adrenoreceptor agonist is indicated, which comprises <u>administering</u> administration of a therapeutically effective amount of a compound of formula (I) according to <u>claim 1</u> any of claims 1 to 10, or a pharmaceutically acceptable salt, solvate, or physiologically functional derivative thereof.

12. (Canceled)

13. (Currently Amended) A pharmaceutical formulation comprising a compound of formula (I) according to <u>claim 1</u> any of claims 1 to 10, or a pharmaceutically acceptable salt, solvate, or physiologically functional derivative thereof, and a pharmaceutically acceptable carrier or excipient, and optionally one or more other therapeutic ingredients.

14. (Canceled)

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15. (Currently Amended) A process for the preparation of a compound of formula (I), according to <u>claim 1</u> any of claims 1 to 10, or a salt, solvate, or physiologically functional derivative thereof, which comprises:

(i) deprotection of deprotecting a protected intermediate, for example of formula (II)

$$Ar^{1a} - CHCH_{2}NR^{23}CR^{1}R^{2}(CH_{2})_{m} - O-(CH_{2})_{p}CR^{1a}R^{2a} - Ar^{2a}$$

$$OR^{24} \qquad (II)$$

or a salt or solvate thereof, wherein R¹, R², R^{1a}, R^{2a}, m, n, p and ____ are as defined for the compound of formula (I), Ar^{1a} represents an optionally protected form of Ar¹; Ar^{2a} represents an optionally protected form of Ar² and R²³ and R²⁴ are each independently either hydrogen or a protecting group, provided that the compound of formula (II) contains at least one protecting group;

(ii) alkylation of an amine of formula

$$Ar^{1a}$$
___CH __CH $_2$ NH R^{23} | OR 24 (VIII)

wherein Ar^{1a}, R²³ and R²⁴ are as defined for formula (II) with a compound of formula (XV):

$$LCR^{1}R^{2}(CH_{2})_{m} CR^{1a}R^{2a}Ar^{2a}$$

$$(XV)$$

wherein ____, Ar², R¹, R², R^{1a}, R^{2a}, m, n and p are as defined for the compound of formula (II) and L is a leaving group as defined for formula (IX);

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wherein said process is optionally followed by one or more of the following steps in any order selected from the group consisting of:

- (i) optional removal of removing any protecting groups;
- (ii) optional separation of separating an enantiomer from a mixture of enantiomers; and
- (iii) optional conversion of converting the product to a corresponding salt, solvate, or physiologically functional derivative thereof.
- 16. (New) A compound of formula (I) as defined in claim
 1, or a salt, solvate or physiologically functional derivative thereof, wherein R¹¹ is –XNR¹⁴C(O)NR¹⁵R¹⁶, and wherein R¹⁴ and R¹⁵ form a 5-, 6-, or 7-membered ring.
- 17. (New) A compound of formula (I) as defined in claim 16, or a salt, solvate or physiologically functional derivative thereof, wherein the 5-, 6-, or 7-membered ring is an imidazolidine ring.
- 18. (New) A compound of formula (I) as defined in claim 17, or a salt, solvate or physiologically functional derivative thereof, wherein the imidazolidine ring is imidazolidine-2,4-dione.
- 19. (New) A compound of formula (I) as defined in claim 1, or a salt, solvate or physiologically functional derivative thereof, where R¹¹ is –XNR¹⁴C(O)OR¹⁵, and wherein R¹⁴ and R¹⁵ form a 5-, 6-, or 7-membered ring.
- 20. (New) A compound of formula (I) as defined in claim 19, or a salt, solvate or physiologically functional derivative thereof, wherein the 5-, 6-, or 7- membered ring is an oxazolidine ring.

- 21. (New) A compound of formula (I) as defined in claim 20, or a salt, solvate or physiologically functional derivative thereof, wherein the oxazolidine ring is oxazolidine-2,4-dione.
- 22. (New) A method according to claim 11, wherein the mammal is a human.
- 23. (New) A process for the preparation of a compound of formula (I), according to claim 1, or a salt, solvate, or physiologically functional derivative thereof, which comprises:

alkylating an amine of formula

$$Ar^{1a}$$
 CH $CH_{2}NHR^{23}$ CH^{24} $(VIII)$

wherein Ar^{1a}, R²³ and R²⁴ are as defined for formula (II) with a compound of formula (XV):

$$LCR^{1}R^{2}(CH_{2})_{m}CR^{1a}R^{2a}Ar^{2a}$$

$$(CH_{2})_{n}O(CH_{2})_{p}CR^{1a}R^{2a}Ar^{2a}$$

$$(XV)$$

wherein ____, Ar², R¹, R², R^{1a}, R^{2a}, m, n and p are as defined for the compound of formula (II) and L is a leaving group as defined for formula (IX);

wherein said process is optionally followed by one or more of the following steps in any order selected from the group consisting of:

- (i) removing any protecting groups;
- (ii) separating an enantiomer from a mixture of enantiomers; and
- (iii) converting the product to a corresponding salt, solvate, or physiologically functional derivative thereof.